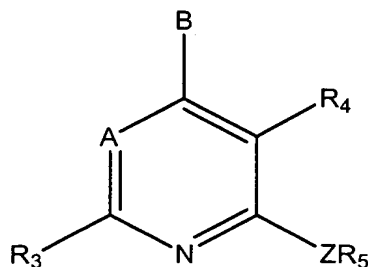


Claim 1 (previously presented). A compound of the formula



or a pharmaceutically acceptable salt thereof, wherein

A is N;

B is  $-NR_1R_2$ ,  $-CR_1R_2R_{11}$ ,  $-C(=CR_2R_{12})R_1$ ,  $-NHCHR_1R_2$ ,  $-OCHR_1R_2$ ,  $-SCHR_1R_2$ ,  $-CHR_2OR_1$ ,  $-CHR_1OR_2$ ,  $-CHR_2SR_1$ ,  $-C(S)R_2$ ,  $-C(O)R_2$ ,  $-CHR_2NR_1R_2$ ,  $-CHR_1NHR_2$ ,  $-CHR_1N(CH_3)R_2$ , or  $-NR_{12}NR_1R_2$ ;

Z is NH, O, S,  $-N(C_1-C_2 \text{ alkyl})-$ ,  $N(C(O)CF_3)-$ , or  $-C(R_{13}R_{14})-$ , wherein  $R_{13}$  and  $R_{14}$  are each, independently, hydrogen, trifluoromethyl or methyl, or one of  $R_{13}$  and  $R_{14}$  is cyano and the other is hydrogen or methyl, or  $-C(R_{13}R_{14})$  is a cyclopropyl group, or Z is nitrogen or CH and forms a five or six membered heterocyclic ring fused with  $R_5$ , which ring optionally comprises two or three further hetero members selected independently from oxygen, nitrogen,  $NR_{12}$ , and  $S(O)_m$ , and optionally comprises from one to three double bonds, and is optionally substituted with halo,  $C_1-C_4$  alkyl,  $-O(C_1-C_4 \text{ alkyl})$ ,  $NH_2$ ,  $NHCH_3$ ,  $N(CH_3)_2$ ,  $CF_3$ , or  $OCF_3$ , with the proviso that said ring does not contain any  $-S-S-$ ,  $-S-O-$ ,  $-N-S-$ , or  $-O-O-$  bonds, and does not comprise more than two oxygen or  $S(O)_m$  heterologous members;

$R_1$  is  $C(O)H$ ,  $C(O)(C_1-C_6 \text{ hydrocarbyl})$ ,  $C(O)(C_1-C_6 [ [$

$\text{[hydrocarbylene)(C}_3\text{-C}_8\text{ cyclohydrocarbyl), C(O)(C}_3\text{-C}_8\text{ cyclohydrocarbylene)}$   
 $\text{)(C}_3\text{-C}_8\text{ cyclohydrocarbyl), C(O)(C}_1\text{-C}_6\text{ hydrocarbylene)(C}_4\text{-C}_8$   
 $\text{heterocyclohydrocarbyl), -C(O)(C}_3\text{-C}_8\text{ cyclohydrocarbylene)(C}_4\text{-C}_8$   
 $\text{heterocyclohydrocarbyl), C}_1\text{-C}_6\text{ hydrocarbyl, C}_3\text{-C}_8\text{ cyclohydrocarbyl, C}_4\text{-C}_8$   
 $\text{heterocyclohydrocarbyl, -(C}_1\text{-C}_6\text{ hydrocarbylene (C}_3\text{-C}_8\text{ cyclohydrocarbyl), C}_3\text{-C}_8$   
 $\text{cyclohydrocarbylene)(C}_3\text{-C}_8\text{ cyclohydrocarbyl), -(C}_1\text{-C}_6\text{ hydrocarbylene)(C}_4\text{-C}_8$   
 $\text{heterocyclohydrocarbyl), -(C}_3\text{-C}_8\text{ cyclohydrocarbylene)(C}_4\text{-C}_8$   
 $\text{heterocyclohydrocarbyl), or -O-aryl, or -O-(C}_1\text{-C}_6\text{ hydrocarbylene)-aryl; wherein said}$   
 $\text{aryl, C}_4\text{-C}_8\text{ heterocyclohydrocarbyl, C}_1\text{-C}_6\text{ hydrocarbyl, C}_3\text{-C}_8\text{ cyclohydrocarbyl,}$   
 $\text{C}_3\text{-C}_8\text{ cyclohydrocarbylene, and C}_1\text{-C}_6\text{ hydrocarbylene groups may each}$   
 $\text{independently be optionally substituted with from one to six fluoro and may each}$   
 $\text{independently be optionally substituted with one or two substituents R}_8\text{ independently}$   
 $\text{selected from the group consisting of C}_1\text{-C}_4\text{ hydrocarbyl, -C}_3\text{-C}_8\text{ cyclohydrocarbyl,}$   
 $\text{hydroxy, chloro, bromo, iodo, CF}_3\text{, -O-(C}_1\text{-C}_6\text{ hydrocarbyl), -O-(C}_3\text{-C}_5$   
 $\text{cyclohydrocarbyl), -O-CO-(C}_1\text{-C}_4\text{ hydrocarbyl), -O-CO-NH(C}_1\text{-C}_4\text{ hydrocarbyl),}$   
 $\text{-O-CO-N(R}_{24}\text{)(R}_{25}\text{), -N(R}_{24}\text{)(R}_{25}\text{), -S(C}_1\text{-C}_4\text{ hydrocarbyl), -S(C}_3\text{-C}_5\text{ cyclohydrocarbyl -}$   
 $\text{-N(C}_1\text{-C}_4\text{ hydrocarbyl)CO(C}_1\text{-C}_4\text{ hydrocarbyl), -NHCO(C}_1\text{-C}_4\text{ hydrocarbyl),}$   
 $\text{-COO(C}_1\text{-C}_4\text{ hydrocarbyl), -CONH(C}_1\text{-C}_4\text{ hydrocarbyl), -CONC}_1\text{-C}_4$   
 $\text{hydrocarbyl)(C}_1\text{-C}_2\text{ hydrocarbyl), CN, NO}_2\text{, -OSO}_2\text{(C}_1\text{-C}_4\text{ hydrocarbyl), S}^+\text{(C}_1\text{-C}_6$   
 $\text{hydrocarbyl)(C}_1\text{-C}_2\text{ hydrocarbyl) I}^-, \text{-SO(C}_1\text{-C}_4\text{ hydrocarbyl) and -SO}_2\text{(C}_1\text{-C}_4$   
 $\text{hydrocarbyl); and wherein the C}_1\text{-C}_6\text{ hydrocarbyl, C}_1\text{-C}_6\text{ hydrocarbylene, C}_5\text{-C}_8$   
 $\text{cyclohydrocarbyl, C}_5\text{-C}_8\text{ cyclohydrocarbylene, and C}_5\text{-C}_8\text{ heterocyclohydrocarbyl}$   
 $\text{moieties of R}_1\text{ may optionally independently contain from one to three double or}$

triple bonds; and wherein the C<sub>1</sub>-C<sub>4</sub> hydrocarbyl moieties and C<sub>1</sub>-C<sub>6</sub> hydrocarbyl moieties of R<sub>8</sub> can optionally independently be substituted with hydroxy, amino, C<sub>1</sub>-C<sub>4</sub> alkyl, aryl, -CH<sub>2</sub>-aryl, C<sub>3</sub>-C<sub>5</sub> cycloalkyl, or -O-(C<sub>1</sub>-C<sub>4</sub> alkyl), and can optionally independently be substituted with from one to six fluoro, and can optionally contain one or two double or triple bonds; and wherein each heterocyclohydrocarbyl group of R<sub>1</sub> contains from one to three heteromoieties selected from oxygen, S(O)<sub>m</sub>, nitrogen, and NR<sub>12</sub>;

R<sub>2</sub> is hydrogen, C<sub>1</sub>-C<sub>12</sub> hydrocarbyl, C<sub>3</sub>-C<sub>8</sub> cyclohydrocarbyl, C<sub>4</sub>-C<sub>8</sub> heterocyclohydrocarbyl, -(C<sub>1</sub>-C<sub>6</sub> hydrocarbylene)(C<sub>3</sub>-C<sub>8</sub> cyclohydrocarbyl), -(C<sub>3</sub>-C<sub>8</sub> cyclohydrocarbylene)(C<sub>3</sub>-C<sub>8</sub> cyclohydrocarbyl), -(C<sub>1</sub>-C<sub>6</sub> hydrocarbylene)(C<sub>4</sub>-C<sub>8</sub> heterocyclohydrocarbyl), -(C<sub>3</sub>-C<sub>6</sub> cyclohydrocarbylene)(C<sub>4</sub>-C<sub>8</sub> heterocyclohydrocarbyl), aryl, -(C<sub>1</sub>-C<sub>6</sub> hydrocarbylene)aryl, or -(C<sub>3</sub>-C<sub>8</sub> cyclohydrocarbylene)(aryl); wherein each of the foregoing R<sub>2</sub> groups may optionally be substituted with from one to three substituents independently selected from chloro, fluoro, and C<sub>1</sub>-C<sub>6</sub> alkyl, wherein one of said one to three substituents can further be selected from bromo, iodo, C<sub>1</sub>-C<sub>6</sub> alkoxy, -OH, -O-CO-(C<sub>1</sub>-C<sub>6</sub> alkyl), -O-CO-N(C<sub>1</sub>-C<sub>4</sub> alkyl)(C<sub>1</sub>-C<sub>2</sub> alkyl), -S(C<sub>1</sub>-C<sub>6</sub> alkyl), -S(O)(C<sub>1</sub>-C<sub>6</sub> alkyl), -S(O)<sub>2</sub>(C<sub>1</sub>-C<sub>6</sub> alkyl), S<sup>+</sup>(C<sub>1</sub>-C<sub>6</sub> alkyl)(C<sub>1</sub>-C<sub>2</sub> alkyl) [I<sup>+</sup>], I<sup>-</sup>, CN, and NO<sub>2</sub>; and wherein the C<sub>1</sub>-C<sub>12</sub> hydrocarbyl, -(C<sub>1</sub>-C<sub>6</sub> hydrocarbylene), and cyclohydrocarbyl groups of 5 - 8 carbon atoms, cyclohydrocarbylene groups of 5 to 8 carbon atoms and heterocyclohydrocarbyl groups of 5 to 8 atoms of R<sub>2</sub> may optionally independently contain from one to three double or triple bonds; and wherein each heterocyclohydrocarbyl group of R<sub>2</sub> contains

from one to three heteromoieties selected from oxygen,  $S(O)_m$ , nitrogen, and  $NR_{12}$ ;

or when  $R_1$  and  $R_2$  are as in  $-NHCHR_1R_2$ ,  $-OCHR_1R_2$ ,  $-SCHR_1R_2$ ,  $-CHR_1R_2$  or  $-NR_1R_2$ ,

$R_1$  and  $R_2$  of B may form a saturated 5- to 8-membered ring which may optionally contain one or

two double bonds and in which one or two of the ring carbons may optionally be replaced by an

oxygen,  $S(O)_m$ , nitrogen or  $NR_{12}$ ; and which carbocyclic ring can optionally be substituted with

from 1 to 3 substituents selected from the group consisting of hydroxy,  $C_1$ - $C_4$  alkyl, fluoro, chloro, bromo, iodo,  $CF_3$ ,  $-O-(C_1-C_4 \text{ alkyl})$ ,  $-O-CO-(C_1-C_4 \text{ alkyl})$ ,

$-O-CO-NH(C_1-C_4 \text{ alkyl})$ ,  $-O-CO-N(C_1-C_4 \text{ alkyl})(C_1-C_2 \text{ alkyl})$ ,  $-NH(C_1-C_4 \text{ alkyl})$ ,

$-N(C_1-C_2 \text{ alkyl})(C_1-C_4 \text{ alkyl})$ ,  $-S(C_1-C_4 \text{ alkyl})$ ,  $-N(C_1-C_4 \text{ alkyl})CO(C_1-C_4 \text{ alkyl})$ ,

$-NHCO(C_1-C_4 \text{ alkyl})$ ,  $-COO(C_1-C_4 \text{ alkyl})$ ,  $-CONH(C_1-C_4 \text{ alkyl})$ ,  $-CON(C_1-C_4$

$\text{alkyl})(C_1-C_2 \text{ alkyl})$ ,  $CN$ ,  $NO_2$ ,  $-OSO_2(C_1-C_4 \text{ alkyl})$ ,  $-SO(C_1-C_4 \text{ alkyl})$ , and  $-SO(C_1-C_4$

$\text{alkyl})$ , wherein one of said one to three substituents can further be selected from phenyl;

$R_3$  is methyl, ethyl, fluoro, chloro, bromo, iodo, cyano, methoxy,  $OCF_3$ ,  $NH_2$ ,  $NH(C_1-C_2 \text{ alkyl})$ ,  $N(CH_3)_2$ ,  $-NHCOCF_3$ ,  $-NHCH_2CF_3$ ,  $S(O)_m(C_1-C_4 \text{ alkyl})$ ,  $CONH_2$ ,  $-CONHCH_3$ ,  $CON(CH_3)_2$ ,  $-CF_3$ , or  $CH_2OCH_3$ ;

$R_4$  is hydrogen,  $C_1$ - $C_4$  hydrocarbyl,  $C_3$ - $C_5$  cycloalkyl,  $-(C_1-C_4 \text{ hydrocarbylene})(C_3-C_5 \text{ cycloalkyl})$ ,  $-(C_3-C_5 \text{ cycloalkylene})(C_3-C_6 \text{ cycloalkyl})$ , cyano, fluoro, chloro, bromo, iodo,  $-OR_{24}$   $C_1$ - $C_6$  alkoxy,  $-O-$  cycloalkyl,  $-O-(C_1-C_4$

hydrocarbylene)(C<sub>3</sub>-C<sub>5</sub> cycloalkyl), -O-(C<sub>3</sub>-C<sub>5</sub> cycloalkylene)(C<sub>3</sub>-C<sub>5</sub> cycloalkyl), -CH<sub>2</sub>SC(S)O(C<sub>1</sub>-C<sub>4</sub> alkyl), CH<sub>2</sub>OCF<sub>3</sub>, CF<sub>3</sub>, amino, nitro, -NR<sub>24</sub>R<sub>25</sub>, -(C<sub>1</sub>-C<sub>4</sub> hydrocarbylene)-OR<sub>24</sub>, -(C<sub>1</sub>-C<sub>4</sub> hydrocarbylene)Cl, -(C<sub>1</sub>-C<sub>4</sub> hydrocarbylene)NR<sub>24</sub>R<sub>25</sub>, -NHCOR<sub>24</sub>, -NHCONR<sub>24</sub>R<sub>25</sub>, -CH=NOR<sub>24</sub>, -NHNOR<sub>24</sub>R<sub>25</sub>, -S(O)<sub>m</sub>R<sub>24</sub>, -C(O)R<sub>24</sub>, -OC(O)R<sub>24</sub>, -C(O)CN, -C(O)NR<sub>24</sub>R<sub>25</sub>, -C(O)NHNOR<sub>24</sub>R<sub>25</sub>, and -COOR<sub>24</sub>, wherein the hydrocarbyl and hydrocarbylene groups of R<sub>4</sub> may optionally independently contain one or two double or triple bonds and may optionally independently be substituted with one or two substituents R<sub>10</sub> independently selected from hydroxy, amino, -NHCOCH<sub>3</sub>, -NHCOCH<sub>2</sub>Cl, -NH(C<sub>1</sub>-C<sub>2</sub> alkyl), -N(C<sub>1</sub>-C<sub>2</sub> alkyl)(C<sub>1</sub>-C<sub>2</sub>alkyl), -COO(C<sub>1</sub>-C<sub>4</sub> alkyl), -COOH, -CO(C<sub>1</sub>-C<sub>4</sub> alkyl), C<sub>1</sub>-C<sub>6</sub> alkoxy, C<sub>1</sub>-C<sub>3</sub> thioalkyl, cyano and nitro, and with one to four substituents independently selected from fluoro and chloro;

R<sub>5</sub> is aryl or heteroaryl and is substituted with from one to four substituents R<sub>27</sub> independently selected from halo, C<sub>1</sub>-C<sub>10</sub> hydrocarbyl, -(C<sub>1</sub>-C<sub>4</sub> hydrocarbylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -(C<sub>1</sub>-C<sub>4</sub> hydrocarbylene)(C<sub>4</sub>-C<sub>8</sub> heterocycloalkyl), -(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -(C<sub>4</sub>-C<sub>8</sub> heterocycloalkyl), -(C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -(C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>4</sub>-C<sub>8</sub> heterocycloalkyl), C<sub>1</sub>-C<sub>4</sub> haloalkyl, C<sub>1</sub>-C<sub>4</sub> haloalkoxy, nitro, cyano, -NR<sub>24</sub>R<sub>25</sub>, -NR<sub>24</sub>COR<sub>25</sub>, -NR<sub>24</sub>CO<sub>2</sub>R<sub>26</sub>, -COR<sub>24</sub>, -OR<sub>25</sub>, -CONR<sub>24</sub>R<sub>25</sub>, -CON(OR<sub>22</sub>)R<sub>23</sub>, -CO<sub>2</sub>R<sub>26</sub>, -C=N(OR<sub>22</sub>)R<sub>23</sub>, and -S(O)<sub>m</sub>R<sub>23</sub>; wherein said C<sub>1</sub>-C<sub>10</sub> alkyl, C<sub>3</sub>-C<sub>8</sub> cycloalkyl, (C<sub>1</sub>-C<sub>4</sub> hydrocarbylene), (C<sub>3</sub>-C<sub>8</sub> cycloalkyl), (C<sub>3</sub>-C<sub>8</sub> cycloalkylene), and (C<sub>4</sub>-C<sub>8</sub> heterocycloalkyl) groups can be optionally substituted with from one to three substituents independently selected from C<sub>1</sub>-C<sub>4</sub> alkyl, C<sub>3</sub>-C<sub>8</sub> cycloalkyl, (C<sub>1</sub>-C<sub>4</sub> hydrocarbylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -(C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>3</sub>-C<sub>8</sub>

cycloalkyl), C<sub>1</sub>-C<sub>4</sub> haloalkyl, hydroxy, C<sub>1</sub>-C<sub>6</sub> alkoxy, nitro, halo, cyano, -NR<sub>24</sub>R<sub>25</sub>, -NR<sub>24</sub>COR<sub>25</sub>, NR<sub>24</sub>CO<sub>2</sub>R<sub>26</sub>, -COR<sub>24</sub>, -OR<sub>25</sub>, -CONR<sub>24</sub>R<sub>25</sub>, CO<sub>2</sub>R<sub>26</sub>, -CO(NOR<sub>22</sub>)R<sub>25</sub>, and -S(O)<sub>m</sub>R<sub>23</sub>; and wherein two adjacent substituents of the R<sub>5</sub> group can optionally form a 5-7 membered ring, saturated or unsaturated, fused to R<sub>5</sub>, which ring optionally can contain one, two, or three heterologous members independently selected from O, S(O)<sub>m</sub>, and N, but not any -S-S-, -O-O-, -S-O-, or -N-S- bonds, and which ring is optionally substituted with C<sub>1</sub>-C<sub>4</sub> alkyl, C<sub>3</sub>-C<sub>8</sub> cycloalkyl, -(C<sub>1</sub>-C<sub>4</sub> alkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -(C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), C<sub>1</sub>-C<sub>4</sub> haloalkyl, nitro, halo, cyano -NR<sub>24</sub>R<sub>25</sub>, NR<sub>24</sub>COR<sub>25</sub>, NR<sub>24</sub>CO<sub>2</sub>R<sub>26</sub>, -COR<sub>24</sub>, -OR<sub>25</sub>, -CONR<sub>24</sub>R<sub>25</sub>, CO<sub>2</sub>R<sub>26</sub>, -CO(NOR<sub>26</sub>)R<sub>25</sub>, or -S(O)<sub>m</sub>R<sub>23</sub>; wherein one of said one to four optional substituents R<sub>27</sub>, can further be selected from -SO<sub>2</sub>NH(C<sub>1</sub>-C<sub>4</sub> alkyl), -SO<sub>2</sub>NH(C<sub>1</sub>-C<sub>4</sub> alkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), SO<sub>2</sub>NH(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -SO<sub>2</sub>NH(C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -SO<sub>2</sub>N(C<sub>1</sub>-C<sub>4</sub> alkyl)(C<sub>1</sub>-C<sub>2</sub> alkyl), -SO<sub>2</sub>NH<sub>2</sub>, -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>4</sub> alkyl), -NHSO<sub>2</sub>(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>4</sub> alkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), and -NHSO<sub>2</sub>(C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl); and wherein the hydrocarbyl, and hydrocarbylene groups of R<sub>5</sub> may independently optionally contain one double or triple bond;

R<sub>6</sub> is hydrogen, C<sub>1</sub>-C<sub>6</sub> alkyl, C<sub>3</sub>-C<sub>8</sub> cycloalkyl, -(C<sub>1</sub>-C<sub>6</sub> alkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), or -(C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), wherein said alkyl and cycloalkyl may optionally be substituted with one hydroxy, methoxy, ethoxy or fluoro group;

or R<sub>6</sub> and R<sub>4</sub> can together form an oxo (=O) group, or can be connected to form a 3-8 membered carbocyclic ring, optionally containing one to three double bonds, and

optionally containing one, two, or three heterologous ring members selected from O, SO<sub>m</sub>, N, and NR<sub>12</sub>, but not containing any -O-O-, -S-O-, -S-S-, or -N-S- bonds, and further optionally substituted with C<sub>1</sub>-C<sub>4</sub> hydrocarbyl or C<sub>3</sub>-C<sub>6</sub> cycloalkyl, wherein said C<sub>1</sub>-C<sub>4</sub> hydrocarbyl substituent may optionally contain one double or triple bond;

R<sub>11</sub> is hydrogen, hydroxy, fluoro, ethoxy, or methoxy;

R<sub>12</sub> is hydrogen or C<sub>1</sub>-C<sub>4</sub> alkyl;

R<sub>22</sub> is independently at each occurrence selected from hydrogen, C<sub>1</sub>-C<sub>14</sub> alkyl, C<sub>1</sub>-C<sub>14</sub> haloalkyl, C<sub>3</sub>-C<sub>6</sub> alkenyl, C<sub>3</sub>-C<sub>6</sub> alkynyl, C<sub>3</sub>-C<sub>8</sub> cycloalkyl, (C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), and (C<sub>1</sub>-C<sub>4</sub>) alkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl);

R<sub>23</sub> is independently at each occurrence selected from C<sub>1</sub>-C<sub>4</sub> alkyl, C<sub>1</sub>-C<sub>4</sub> haloalkyl, C<sub>2</sub>-C<sub>8</sub> alkoxyalkyl, C<sub>3</sub>-C<sub>8</sub> cycloalkyl, -(C<sub>1</sub>-C<sub>4</sub> alkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -(C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), aryl, -(C<sub>1</sub>-C<sub>4</sub> alkylene)aryl, piperidine, pyrrolidine, piperazine, N-methylpiperazine, morpholine, and thiomorpholine;

R<sub>24</sub> and R<sub>25</sub> are independently at each occurrence selected from hydrogen, -C<sub>1</sub>-C<sub>4</sub> alkyl, C<sub>1</sub>-C<sub>4</sub> haloalkyl, -(C<sub>1</sub>-C<sub>4</sub> alkylene)OH, -(C<sub>1</sub>-C<sub>4</sub> alkylene)-O-(C<sub>1</sub>-C<sub>4</sub> alkyl), -(C<sub>1</sub>-C<sub>4</sub> alkylene)-O-(C<sub>3</sub>-C<sub>5</sub> cycloalkyl), C<sub>3</sub>-C<sub>8</sub> cycloalkyl, -(C<sub>1</sub>-C<sub>4</sub> alkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -(C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>3</sub>-C<sub>8</sub> cycloalkyl), -C<sub>4</sub>-C<sub>8</sub> heterocyclohydrocarbyl, -(C<sub>1</sub>-C<sub>4</sub> alkylene)(C<sub>4</sub>-C<sub>8</sub> heterocyclohydrocarbyl), -(C<sub>3</sub>-C<sub>8</sub> cycloalkylene)(C<sub>4</sub>-C<sub>8</sub> heterocyclohydrocarbyl), aryl, and -(C<sub>1</sub>-C<sub>4</sub> alkylene)(aryl), wherein the -C<sub>4</sub>-C<sub>8</sub> heterocyclohydrocarbyl groups can each independently optionally be substituted with aryl, CH<sub>2</sub>-aryl, or C<sub>1</sub>-C<sub>4</sub> alkyl, and can optionally contain one or two double or triple bonds; or, when R<sub>24</sub> and R<sub>25</sub> are as NR<sub>24</sub>R<sub>25</sub>, -C(O)NR<sub>24</sub>R<sub>25</sub>, -(C<sub>1</sub>-C<sub>4</sub> alkylene)NR<sub>24</sub>R<sub>25</sub>, or -NHCONR<sub>24</sub>R<sub>25</sub>, then NR<sub>24</sub>R<sub>25</sub> may further optionally form a 4 to 8 membered

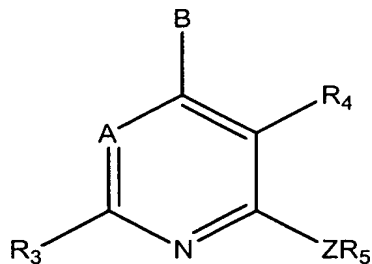
heterocyclic ring optionally containing one or two further hetero members independently selected from  $S(O)_m$ , oxygen, nitrogen, and  $NR_{12}$ , and optionally containing from one to three double bonds;

$R_{26}$  is independently at each occurrence selected from  $C_1$ - $C_4$  alkyl,  $C_1$ - $C_4$  haloalkyl,  $C_3$ - $C_8$  cycloalkyl,  $-(C_1$ - $C_4$  alkylene)( $C_3$ - $C_8$  cycloalkyl),  $-(C_3$ - $C_8$  cycloalkylene)( $C_3$ - $C_8$  cycloalkyl), aryl, and  $-(C_1$ - $C_4$  alkylene)(aryl); and

wherein each  $m$  is independently zero, one, or two,

with the proviso that heterocyclohydrocarbylene groups of the compound of formula I, do not comprise any  $-S-S-$ ,  $-S-O-$ ,  $-N-S-$ , or  $-O-O-$  bonds, and do not comprise more than two oxygen or  $S(O)_m$  heterologous members.

Claim 2 (previously presented). A compound according to claim 1 of the formula



or a pharmaceutically acceptable salt thereof, wherein

A is N;

B is  $-NR_1R_2$ ,  $-CR_1R_2R_{11}$ ,  $-C(=CR_2R_{12})R_1$ ,  $-NHCHR_1R_2$ ,  $-OCHR_1R_2$ ,  $-SCHR_1R_2$ ,  $-CHR_2OR_{12}$ ,  $-CHR_2SR_{12}$ ,  $-C(S)R_2$  or  $-C(O)R_2$ ;

Z is  $-NH$ , O, S,  $N(C_1$ - $C_2$  alkyl) or  $C(R_{13}R_{14})$  wherein  $R_{13}$  and  $R_{14}$  are each independently, hydrogen, trifluoromethyl or methyl or one of  $R_{13}$  and  $R_{14}$  is cyano and the other is hydrogen or methyl;



R<sub>1</sub> is C<sub>1</sub>-C<sub>6</sub> hydrocarbyl which may optionally be substituted with one or two substituents R<sub>8</sub> independently selected from the group consisting of hydroxy, fluoro, chloro, bromo, iodo, CF<sub>3</sub>, C<sub>1</sub>-C<sub>4</sub> alkoxy, -O-CO-(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl), -O-CO-NH(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl), -O-CO-N(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl)(C<sub>1</sub>-C<sub>2</sub> hydrocarbyl), -NH(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl), -N(C<sub>1</sub>-C<sub>2</sub> alkyl)(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl), -S(C<sub>1</sub>-C<sub>4</sub> alkyl), -N(C<sub>1</sub>-C<sub>4</sub>)CO(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl), -NHCO(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl), -COO(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl)hydrocarbyl, -CONH(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl), -CON(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl)(C<sub>1</sub>-C<sub>2</sub> alkyl), CN, NO<sub>2</sub>, -SO(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl) and -SO<sub>2</sub>(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl), and wherein said C<sub>1</sub>-C<sub>6</sub> hydrocarbyl and the (C<sub>1</sub>-C<sub>4</sub>)hydrocarbyl moieties in the foregoing R<sub>1</sub> groups may optionally contain one carbon-carbon double or triple bond;

R<sub>2</sub> is C<sub>1</sub>-C<sub>12</sub> hydrocarbyl, aryl or -(C<sub>1</sub>-C<sub>4</sub> hydrocarbylene)aryl wherein said aryl is phenyl, naphthyl, thienyl, benzothienyl, pyridyl, quinolyl, pyrazinyl, pyrimidyl, imidazolyl, furanyl, benzofuranyl, benzothiazolyl, isothiazolyl, benzisothiazolyl, benzisoxazolyl, benzimidazolyl, indolyl, or benzoxazolyl; 3- to 8-membered cycloalkyl or -(C<sub>1</sub>-C<sub>6</sub> alkylene)cycloalkyl, wherein one or two of the ring carbons of said cycloalkyl having at least 4 ring members and the cycloalkyl moiety of said -(C<sub>1</sub>-C<sub>6</sub> alkylene)cycloalkyl having at least 4 ring members may optionally be replaced by an oxygen or sulfur atom or by N-R<sub>9</sub> wherein R<sub>9</sub> is hydrogen or C<sub>1</sub>-C<sub>4</sub> alkyl; and wherein each of the foregoing R<sub>2</sub> groups may optionally be substituted with from one to three substituents independently selected from chloro, fluoro and C<sub>1</sub>-C<sub>4</sub> alkyl, or with one substituent selected from bromo, iodo, C<sub>1</sub>-C<sub>6</sub> alkoxy, -O-CO-(C<sub>1</sub>-C<sub>6</sub> alkyl), -O-CO-N(C<sub>1</sub>-C<sub>4</sub> alkyl)(C<sub>1</sub>-C<sub>2</sub> alkyl), -S(C<sub>1</sub>-C<sub>6</sub> alkyl), CN, NO<sub>2</sub>,

-SO(C<sub>1</sub>-C<sub>4</sub> alkyl), and -SO<sub>2</sub>(C<sub>1</sub>-C<sub>4</sub> alkyl), and wherein said C<sub>1</sub>-C<sub>12</sub> hydrocarbyl and the C<sub>1</sub>-C<sub>4</sub> hydrocarbylene moiety of said -(C<sub>1</sub>-C<sub>4</sub> hydrocarbylene)aryl may optionally contain one carbon-carbon double or triple bond;

or -NR<sub>1</sub>R<sub>2</sub> or -CR<sub>1</sub>R<sub>2</sub>R<sub>11</sub> may form a saturated 5- to 8-membered carbocyclic ring which may optionally contain one or two carbon-carbon double bonds and in which one or two of the ring carbons may optionally be replaced by an oxygen or sulfur atom;

R<sub>3</sub> is methyl, ethyl, fluoro, chloro, bromo, iodo, cyano, methoxy, OCF<sub>3</sub>, methylthio, methylsulfonyl, CH<sub>2</sub>OH, or CH<sub>2</sub>OCH<sub>3</sub>;

R<sub>4</sub> is hydrogen, C<sub>1</sub>-C<sub>4</sub> hydrocarbyl, fluoro, chloro, bromo, iodo, C<sub>1</sub>-C<sub>4</sub> alkoxy, trifluoromethoxy, -CH<sub>2</sub>OCH<sub>3</sub>, -CH<sub>2</sub>OCH<sub>2</sub>CH<sub>3</sub>, -CH<sub>2</sub>CH<sub>2</sub>OCH<sub>3</sub>, -CH<sub>2</sub>OF<sub>3</sub>, CF<sub>3</sub>, amino, nitro, -NH(C<sub>1</sub>-C<sub>4</sub> alkyl), -N(CH<sub>3</sub>)<sub>2</sub>, -NHCOCH<sub>3</sub>, -NHCONHCH<sub>3</sub>, -SO<sub>n</sub>(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl) wherein n is 0, 1 or 2, cyano, hydroxy, -CO(C<sub>1</sub>-C<sub>4</sub> hydrocarbyl), -CHO, cyano or -COO(C<sub>1</sub>-C<sub>4</sub> alkyl) wherein said C<sub>1</sub>-C<sub>4</sub> hydrocarbyl may optionally contain one double or triple bond and may optionally be substituted with one substituent selected from hydroxy, amino, -NHCOCH<sub>3</sub>, -NH(C<sub>1</sub>-C<sub>2</sub> alkyl), -N(C<sub>1</sub>-C<sub>2</sub> alkyl)<sub>2</sub>, -COO(C<sub>1</sub>-C<sub>4</sub> alkyl), -CO(C<sub>1</sub>-C<sub>4</sub> alkyl), C<sub>1</sub>-C<sub>3</sub> alkoxy, C<sub>1</sub>-C<sub>3</sub> thioalkyl, fluoro, chloro, cyano and nitro;

R<sub>5</sub> is phenyl, naphthyl, thienyl, benzothienyl, pyridyl, quinolyl, pyrazinyl, pyrimidyl, furanyl, benzofuranyl, benzothiazolyl, or indolyl, wherein each of the above groups R<sub>5</sub> is substituted with from one to three substituents independently selected from fluoro, chloro, C<sub>1</sub>-C<sub>6</sub> alkyl, and C<sub>1</sub>-C<sub>6</sub> alkoxy, or with one substituent selected from hydroxy, iodo, bromo, formyl, cyano, nitro, trifluoromethyl, amino,

-(C<sub>1</sub>-C<sub>6</sub> alkyl)O(C<sub>1</sub>-C<sub>6</sub>)alkyl, -NHCH<sub>3</sub>, -N(CH<sub>3</sub>)<sub>2</sub>, -COOH, -COO(C<sub>1</sub>-C<sub>4</sub> alkyl), -CO(C<sub>1</sub>-C<sub>4</sub> alkyl), -SO<sub>2</sub>NH(C<sub>1</sub>-C<sub>4</sub> alkyl), -SO<sub>2</sub>N(C<sub>1</sub>-C<sub>4</sub> alkyl)(C<sub>1</sub>-C<sub>2</sub> alkyl), -SO<sub>2</sub>NH<sub>2</sub>, -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>4</sub> alkyl), -S(C<sub>1</sub>-C<sub>6</sub> alkyl) and -SO<sub>2</sub>(C<sub>1</sub>-C<sub>6</sub> alkyl) , and wherein the C<sub>1</sub>-C<sub>4</sub> alkyl and C<sub>1</sub>-C<sub>6</sub> alkyl moieties of the foregoing R<sub>5</sub> groups may optionally be substituted with one or two fluoro groups or with one substituent selected from hydroxy, amino, methylamino, dimethylamino and acetyl;

R<sub>11</sub> is hydrogen, hydroxy, fluoro, or methoxy;

R<sub>12</sub> is hydrogen or C<sub>1</sub>-C<sub>4</sub> alkyl; and

or a pharmaceutically acceptable salt of such compound.

Claim 3 (previously presented) A compound according to claim 2 wherein B is -NR, R<sub>2</sub>, -NHCHR<sub>1</sub>R<sub>2</sub> or -OCHR<sub>1</sub>R<sub>2</sub>; R<sub>1</sub> is C<sub>1</sub>-C<sub>6</sub> hydrocarbyl, which may optionally be substituted with one hydroxy, fluoro, CF<sub>3</sub>, or C<sub>1</sub>-C<sub>2</sub> alkoxy group and may optionally contain one double or triple bond; and R<sub>2</sub> is benzyl or C<sub>1</sub>-C<sub>6</sub> hydrocarbyl which may optionally contain one carbon-carbon double or triple bond, wherein said C<sub>1</sub>-C<sub>6</sub> alkyl or the phenyl moiety of said benzyl may optionally be substituted with fluoro, CF<sub>3</sub>, C<sub>1</sub>-C<sub>2</sub> alkoxy.

Claim 4 (previously presented) A compound according to claim 2 wherein R<sub>1</sub> is C<sub>1</sub>-C<sub>6</sub> hydrocarbyl which may be substituted by fluoro, CF<sub>3</sub>, hydroxy; C<sub>1</sub>-C<sub>2</sub> alkoxy and which may optionally contain one carbon-carbon double or triple bond.

Claim 5 (original) A compound according to claim 2 wherein R<sub>2</sub> is C<sub>1</sub>-C<sub>4</sub> alkyl which

may optionally be substituted by fluoro, chloro, CF<sub>3</sub>, C<sub>1</sub>--C<sub>4</sub> alkoxy.

Claim 6 (original) A compound according to claim 2 wherein R<sub>3</sub> is methyl, chloro, or methoxy.

Claim 7 (original) A compound according to claim 2 wherein R<sub>4</sub> is methyl, -CH<sub>2</sub>OH, cyano, trifluoromethoxy, methoxy, chloro, trifluoromethyl, -COOCH<sub>3</sub>, -CH<sub>2</sub>Cl, -CH<sub>2</sub>F, ethyl, amino or nitro.

Claim 8 (original) A compound according to claim 2 wherein R<sub>5</sub> is phenyl substituted with two or three substituents.

Claim 9 (original) A compound according to claim 2 wherein R<sub>5</sub> is pyridyl substituted with two or three substituents.

Claim 10 (original) A compound according to claim 8 wherein said substituents are selected, independently, from fluoro, chloro bromo, iodo, C<sub>1</sub>-C<sub>4</sub> alkoxy, trifluoromethyl, C<sub>1</sub>-C<sub>6</sub> alkyl which may optionally be substituted with one hydroxy, C<sub>1</sub>-C<sub>4</sub> alkoxy or fluoro group and which may optionally contain one carbon-carbon double or triple bond, -(C<sub>1</sub>-C<sub>4</sub> alkylene)O(C<sub>1</sub>-C<sub>2</sub> alkyl), C<sub>1</sub>-C<sub>3</sub> hydroxyalkyl, hydroxy, formyl, COO(C<sub>1</sub>-C<sub>4</sub> alkyl).

Claim 11 (original) A compound according to claim 9 wherein said substituents are

selected, independently from fluoro, chloro, bromo, iodo, C<sub>1</sub>-C<sub>4</sub> alkoxy, trifluoromethyl, C<sub>1</sub>-C<sub>6</sub> alkyl which may optionally be substituted with one hydroxy, C<sub>1</sub>-C<sub>4</sub> alkoxy or fluoro group and which may optionally contain one carbon-carbon double or triple bond, -(C<sub>1</sub>-C<sub>4</sub> alkylene)O(C<sub>1</sub>-C<sub>2</sub> alkyl), C<sub>1</sub>-C<sub>3</sub> hydroxyalkyl, hydroxy, formyl, -COO(C<sub>1</sub>-C<sub>2</sub> alkyl) -(C<sub>1</sub>-C<sub>2</sub> alkylene)amino, and -(C(O)(C<sub>1</sub>-C<sub>4</sub> alkyl).

Claim 12 (previously presented) A compound according to claim 1, wherein said compound is 4-(1-ethyl-propoxy)-2,5-dimethyl-6-(2,4,6-trimethyl-benzyl)-pyrimidine; [2,5-dimethyl-6-trimethyl-phenoxy)-pyrimidin-4-yl](1-ethyl-propyl)-amine;

(1-ethyl-propyl)-[2-methyl-5-nitro-6-(2,4,6-trimethyl-pyridin-3-yloxy)-pyrimidin-4-yl]-amine;

(N-(1-ethyl-propyl)-2methyl-5nitro-N'-N= (2,4,6-trimethyl-pyridin-3-yl)-pyrimidine-4,6-diamine;

4-(1-ethylpropoxy)-2,5-dimethyl-6-(2,4,6-trimethylphenoxy)-pyrimidine;

N-butyl-N-ethyl-2,5-dimethyl-N'-(2,4,6-trimethylphenyl)-pyrimidine-4,6-diamine; or

6-(1-ethyl-propoxy)-2-methyl-N4-(2,4,6-trimethyl-phenyl)-pyrimidine-4,5-diamine;

or a pharmaceutically acceptable salt of one of the above compounds.

Claim 13 (previously presented). A pharmaceutical composition for the treatment of (a) a disorder or condition the treatment of which can be effected or facilitated by antagonizing CRF or (b) a disorder or condition selected from inflammatory disorders, pain, asthma, psoriasis and allergies; generalized anxiety disorder; panic; phobias;

obsessive-compulsive disorder; post-traumatic stress disorder; sleep disorders induced by stress; pain perception; mood disorders, mood disorders associated with premenstrual syndrome, and postpartum depression; dysthemia; bipolar disorders; cyclothymia; chronic fatigue syndrome; stress-induced headache; cancer; irritable bowel syndrome, Crohn's disease; spastic colon; post operative ileus; ulcer; diarrhea; stress-induced fever; human immunodeficiency virus infections; neurodegenerative diseases, gastrointestinal diseases; eating disorder; hemorrhagic stress; chemical dependencies or addictions; drug or alcohol withdrawal symptoms; stress-induced psychotic episodes; euthyroid sick syndrome; syndrome of inappropriate antidiuretic hormone; obesity; infertility; head trauma; spinal cord trauma; ischemic neuronal damage; excitotoxic neuronal damage; epilepsy; stroke; immune dysfunctions; muscular spasms; urinary incontinence; senile dementia of the Alzheimer's type; multi infarct dementia; amyotrophic lateral sclerosis; hypertension; tachycardia; congestive heart failure; osteoporosis; premature birth; hypoglycemia, and Syndrome X in a mammal or bird, comprising an amount of a compound according to claim 1 that is effective in the treatment of such disorder or condition, and a pharmaceutically acceptable carrier.

Claim 14 (previously presented). A pharmaceutical composition according to claim 13 for the treatment of a disorder selected from inflammatory disorders; pain, asthma, psoriasis and allergies; generalized anxiety disorder; panic; phobias; obsessive compulsive disorder; post-traumatic stress disorder; sleep disorders induced by stress; pain perception; mood disorders dysthemia; bipolar disorders; cyclothymia; fatigue

syndrome; stress induced headache; cancer; irritable bowel syndrome, Crohn's disease; spastic colon; human immunodeficiency virus (HIV) infections; neurodegenerative diseases; gastrointestinal diseases; eating disorders; chemical dependencies and addictions; obesity; infertility; head traumas; spinal cord trauma; ischemic neuronal damage; excitotoxic neuronal damage; epilepsy; stroke; immune dysfunctions; muscular spasms; urinary incontinence; senile dementia of the Alzheimer's type; multi infarct dementia; amyotrophic lateral sclerosis; and hypoglycemia in a mammal.

Claims 15 to 28 (cancelled).

Claim 29 (previously presented). A compound as claimed in claim 1 wherein  $R_{24}$  and  $R_{25}$  are selected from  $-CF_3$ ,  $-CHF_2$ ,  $CF_2CF_3$ , and  $CH_2CF_3$ .

Claim 30 (previously presented) A pharmaceutical composition as claimed in claim 13 for treatment of a mood disorder selected from the group consisting of rheumatoid arthritis and osteoarthritis, pain, asthma, psoriasis and allergies.

Claim 31 (previously presented) A pharmaceutical composition as claimed in claim 13 for treatment of an inflammatory disorder selected from the group consisting of rheumatoid arthritis and osteoarthritis.

Claim 32 (previously presented). A pharmaceutical composition as claimed in claim

14 for treatment of depression, selected from the group consisting of major depression, single episode depression, recurrent depression, and child abuse induced depression.

Claim 33 (previously presented) A pharmaceutical composition as claimed in claim 14 for treatment of neurodegenerative diseases selected from the group consisting of Alzheimer's disease, Parkinson's disease and Huntington's disease.

Claim 34 (previously presented) A pharmaceutical composition as claimed in claim 14 for treatment of chemical dependencies or addictions, selected from the group consisting of dependencies or addictions to alcohol, cocaine, heroin, benzodiazapines, or other drugs.

Claim 35 (previously presented) A pharmaceutical composition as claimed in claim 14 for treatment of cerebral ischemia.

Claim 36 (previously presented). A pharmaceutical composition as claimed in claim 14 for treatment of immune dysfunctions induced by stress selected from the group consisting of porcine stress syndrome, bovine shipping fever, equine paroxysmal fibrillation, confinement dysfunction in chicken, sheering stress in sheep, and human animal interaction stress in dogs.

Claim 37 (previously presented) A pharmaceutical composition as claimed in claim



14 for treatment of fibromyalgia.

Claim 38 (previously presented) A pharmaceutical composition as claimed in claim 14 for treatment of anorexia or bulimia nervosa.

Claim 39 (previously presented). A pharmaceutical composition as claimed in claim 44 for treatment of cerebral ischemia, selected from the group consisting of cerebral hippocampal ischemia.

Claim 40 (previously presented). A pharmaceutical composition as claimed in claim 14 for treatment of social phobia, agoraphobia, or specific phobias.

Claim 41 (previously presented). The pharmaceutical composition according to claim 13 wherein the pain perception is fibromyalgia.

Claim 42 (previously presented). The pharmaceutical composition according to claim 13 wherein the ischemic neuronal damage is cerebral ischemia.

Claim 43 (previously presented). The pharmaceutical composition according to claim 14 for the treatment of depression or postpartum depression.

Claim 44 (previously presented). The pharmaceutical composition according to claim 14 wherein the ischemic neuronal damage is cerebral ischemia.

Claim 45 (previously presented). The pharmaceutical composition according to claim 14 wherein the mammal is a human.

Claim 46 (new) The compound of claim 1, where Z is NH, O, S, NC(OC)CF<sub>3</sub>, or CR<sub>13</sub>R<sub>14</sub>.

Claim 47 (new) The compound of Claim 29, wherein Z is NH, S, NC(OC)CF<sub>3</sub>, or CR<sub>13</sub>R<sub>14</sub>.

Claim 48 (new) The compound of claim 29, wherein Z is NH.